

Electrical and Thermoelectrical Properties of *TlMnS₂* and *TlMnSe₂*

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The synthesis regimes of *TlMnSe₂* and *TlMnS₂* phases have been worked out. X-ray analysis showed that *TlMnS₂* is crystallized in tetragonal structure with elementary cell parameters $a=7.74$; $c=30.60\approx$; roentgen density $\rho_x=6.40$ g/cm³. *TlMnSe₂* is crystallized in hexagonal structure with elementary cell parameters $a=6.53$; $c=23.96\approx$; $z=8$; $\rho_x=6.71$ g/cm³.

The electric and thermoelectric properties of *TlMnSe₂* and *TlMnS₂* were studied. In the temperature range 130 - 315 K, temperature dependence of the conductivity (σ) of *TlMnS₂* increases exponentially with increasing temperature, i.e. $\sigma(T)$ -dependence had a semiconductor nature. It was shown that $\sigma(T)$ -dependence of *TlMnS₂* consists of three various regions with following activation energies: 0.178, 0.08 and 0.44eV.

The temperature dependence of the conductivity of *TlMnSe₂* had a metallic nature. Temperature dependence of the thermoelectromotive force (α) in *TlMnSe₂* was studied. The thermo-e.m.f. sign corresponded to the p-type conductivity of *TlMnSe₂* in the temperature range 88 - 300K. With increasing temperature from 88 to 300K, the value of the thermo- e.m.f. in *TlMnSe₂* increased from 77 to 200 μ V/K. At $T=194$ K an anomaly was revealed in the dependence $\alpha(T)$.

The low-temperature branch of $\alpha(T)$ -dependence in *TlMnSe₂* had a linear character with extrapolation to $T=0$ according to metallic formula for thermo-e.m.f.:

$$\alpha(T) = \frac{\pi^2}{3} \cdot \frac{k^2 T}{e} \left(\frac{\partial \ln \sigma}{\partial E} \right)_{E=E_F},$$

where k is Boltzmann constant; e -electron charge; E_F -Fermi energy and $kT \ll E_F$.